

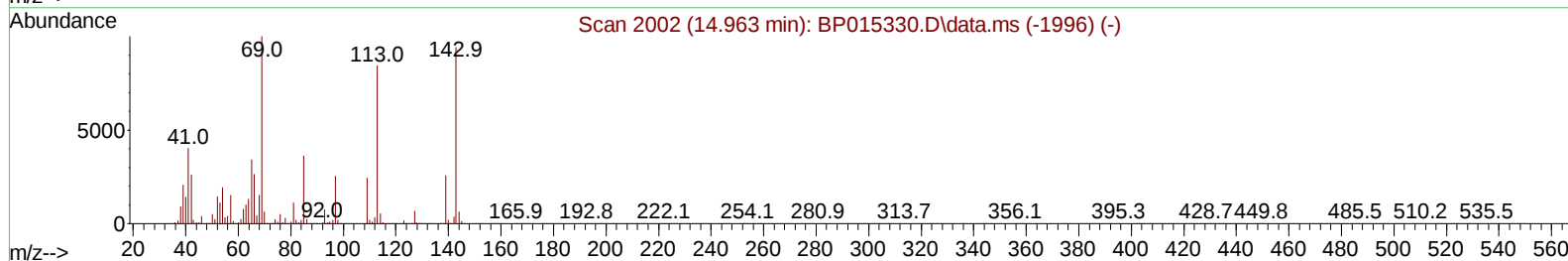
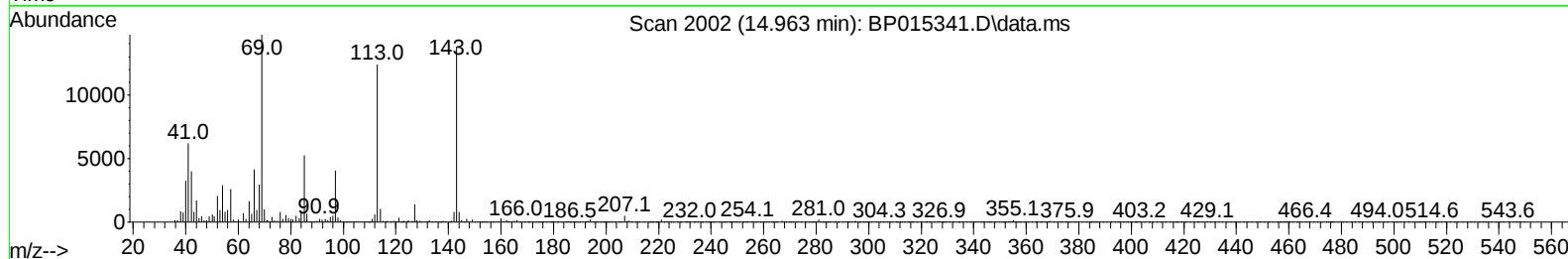
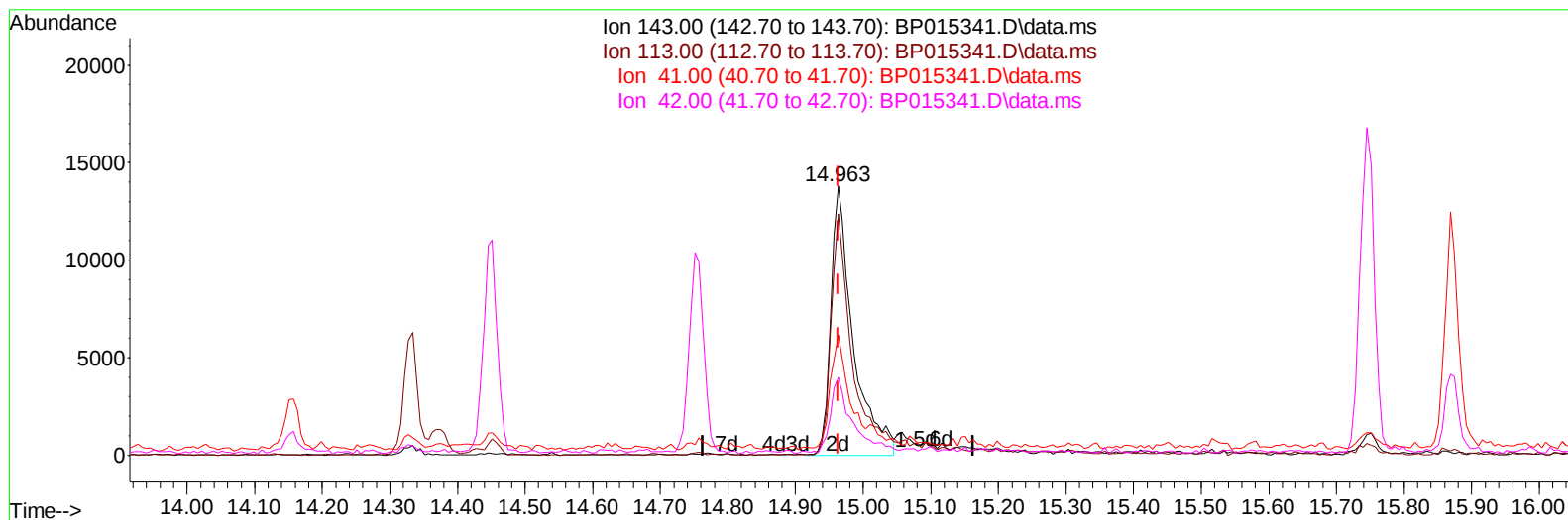
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP053023\
Data File : BP015341.D
Acq On : 30 May 2023 18:49
Operator : CG/JU
Sample : 02961-09
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
H0FQ1

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 05/31/2023
Supervised By :Jagrut Upadhyay 05/31/2023

Quant Time: May 31 00:46:02 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 30 15:23:22 2023
Response via : Initial Calibration



TIC: BP015341.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.963min (-0.000) 4.58 ng/u1 m

response 31601

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	67.30	89.82#
41.00	36.10	44.76#
42.00	24.50	28.84

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Quant Time: May 31 01:09:38 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051123.MA.M
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.110	152	150598	20.000	ng/ul	0.00
20) Naphthalene-d8	10.940	136	667665	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.757	164	470368	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.510	188	1115563	20.000	ng/ul	0.00
79) Chrysene-d12	21.586	240	1114474	20.000	ng/ul	-0.01
88) Perylene-d12	24.151	264	1232819	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.416	96	16213	4.183	ng/uL	0.00
4) Pyridine-d5	3.869	84	59858	5.448	ng/ul	0.00
7) Phenol-d5	7.263	99	66817	4.779	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.428	67	218454	27.051	ng/ul	0.00
11) 2-Chlorophenol-d4	7.640	132	206313	20.068	ng/ul	0.00
15) 4-Methylphenol-d8	8.822	113	131957	11.843	ng/ul	-0.01
21) Nitrobenzene-d5	9.287	128	152550	29.077	ng/ul	0.00
24) 2-Nitrophenol-d4	10.016	143	161762	27.710	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.557	165	249709	23.705	ng/ul	0.00
31) 4-Chloroaniline-d4	11.081	131	247544	15.883	ng/ul	0.00
46) Dimethylphthalate-d6	14.157	166	1139470	32.058	ng/ul	0.00
49) Acenaphthylene-d8	14.451	160	1209990	29.765	ng/ul	0.00
54) 4-Nitrophenol-d4	14.963	143	31601m	4.579	ng/ul	0.00
60) Fluorene-d10	15.745	176	947269	31.584	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.869	200	206807	29.416	ng/ul	-0.01
73) Anthracene-d10	17.610	188	1617021	31.655	ng/ul	0.00
81) Pyrene-d10	19.827	212	2046304	31.658	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.980	264	2051675	33.525	ng/ul	-0.01

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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